

Data Path : Z:\svoasrv\HPCHEM1\BNA_M\Data\BM011422\
 Data File : BM033740.D
 Acq On : 14 Jan 2022 22:23
 Operator : CG/JU
 Sample : N1144-01
 Misc :
 ALS Vial : 17 Sample Multiplier: 1

Instrument :
 BNA_M
 ClientSampleId :
 RO-ETGI-352

Quant Time: Jan 17 00:28:07 2022
 Quant Method : Z:\svoasrv\HPCHEM1\BNA_M\Methods\8270-BM010522.M
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION
 QLast Update : Thu Jan 13 07:49:01 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) 1,4-Dichlorobenzene-d4	7.960	152	166977	20.000	ng	-0.01
21) Naphthalene-d8	10.778	136	713726	20.000	ng	#-0.01
39) Acenaphthene-d10	14.601	164	464486	20.000	ng	0.00
64) Phenanthrene-d10	17.336	188	959183	20.000	ng	0.00
76) Chrysene-d12	21.501	240	624120	20.000	ng	0.00
86) Perylene-d12	23.912	264	679312	20.000	ng	-0.02
System Monitoring Compounds						
5) 2-Fluorophenol	5.507	112	1188818	112.287	ng	0.00
7) Phenol-d6	7.119	99	1551121	108.704	ng	0.00
23) Nitrobenzene-d5	9.125	82	1020460	66.030	ng	-0.01
42) 2,4,6-Tribromophenol	16.083	330	647743	133.516	ng	0.00
45) 2-Fluorobiphenyl	13.230	172	2218556	62.560	ng	-0.01
79) Terphenyl-d14	19.948	244	2634191	75.971	ng	-0.01
Target Compounds						
						Qvalue
32) Benzoic acid	10.037	122	96553	13.274	ng	# 86
71) Phenanthrene	17.377	178	265048	4.829	ng	98
75) Fluoranthene	19.383	202	488889	8.485	ng	98
78) Pyrene	19.742	202	444518	10.116	ng	98
81) Benzo(a)anthracene	21.483	228	185629	4.386	ng	96
83) Chrysene	21.536	228	186685	4.609	ng	96
84) Bis(2-ethylhexyl)phtha...	21.412	149	158419	5.299	ng	99
87) Indeno(1,2,3-cd)pyrene	26.436	276	136625	2.727	ng	96
88) Benzo(b)fluoranthene	23.177	252	279334	6.450	ng	# 91
90) Benzo(a)pyrene	23.806	252	181562	5.051	ng	# 95
92) Benzo(g,h,i)perylene	27.212	276	135391	3.310	ng	96

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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